

Comment on “Quasiparticle Anisotropy and Pseudogap Formation from the Weak-Coupling Renormalization Group Point of View”

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In their recent Letter, Katanin and Kampf [1] (KK) reported numerical results for the self-energy $\Sigma(\mathbf{k}; \varepsilon)$, $\varepsilon \in \mathbb{R}$, of the single-band Hubbard Hamiltonian (hereafter ‘single-band’ will be implicit) in two space dimensions (i.e. $d = 2$), obtained through employing the functional renormalization-group (fRG) formalism (see references in [1], and [2]) at the one-loop level. Several of the results by KK are in full conformity with the *exact* formal results reported earlier in [3, 4, 5]. This, as we shall elaborate below, strengthens one’s confidence in the reliability of the fRG in dealing with models of strongly-correlated fermions.

(i) *The quasi-particle (qp) weight Z_F at the Fermi surface.* — Amongst other things, KK showed that [1] “At van Hove (vH) band fillings and at low temperatures, the quasiparticle weight along the Fermi surface (FS) continuously vanishes on approaching the $(\pi, 0)$ point” In [3] we have obtained a general expression for the momentum-distribution function $n(\mathbf{k})$ at $\mathbf{k} = \mathbf{k}_F^\mp$, i.e. infinitesimally in- and outside Fermi sea (here we suppress the spin indices encountered in [3, 4, 5]), which on the basis of the available numerical results at the time we have shown to reduce to (Eq. (99) in [3]): $n(\mathbf{k}_F^\mp) = (a + Ub^\mp)/(a + 2Ub^\mp)$, where a and b^\mp are components of vectors $\mathbf{a}(\mathbf{k}_F) \equiv \nabla_{\mathbf{k}} \varepsilon_{\mathbf{k}}|_{\mathbf{k}=\mathbf{k}_F}$ and $\mathbf{b}(\mathbf{k}_F^\mp)$ along the outward unit vector normal to Fermi surface \mathcal{S}_F at $\mathbf{k} = \mathbf{k}_F$; here $\varepsilon_{\mathbf{k}}$ is the non-interacting energy dispersion and $\mathbf{b}(\mathbf{k})$ the gradient of a well-defined scalar ground-state (GS) correlation function. Stability of the latter GS is tantamount to the satisfaction of [3] $b^- > a/(U\Lambda^-)$, $b^+ < -a/U$, where $\Lambda^- \equiv n(\mathbf{k}_F^-)/[1 - n(\mathbf{k}_F^-)] > 0$. Our above expression for $n(\mathbf{k}_F^\mp)$ makes explicit that (a) $n(\mathbf{k}_F^\mp) \gtrless 1/2$, (b) $n(\mathbf{k}_F^\mp) \rightarrow 1/2$, i.e. $Z_F \equiv n(\mathbf{k}_F^-) - n(\mathbf{k}_F^+) \rightarrow 0$, for $Ub^\mp \rightarrow \infty$, and (c) $n(\mathbf{k}_F^\mp) = 1/2$, i.e. $Z_F = 0$, for $a = 0$, that is for \mathbf{k}_F at vH points. In Fig. 1 we compare our results for Z_F with those determined by KK [1], *disregarding* the dependence of b^\mp on the direction of \mathbf{k}_F .

(ii) *The single-particle spectral function $A(\mathbf{k}; \varepsilon)$ for \mathbf{k} in the pseudogap (PG) region.* — KK observed that [1] “The qp weight suppression [for $\mathbf{k}_F \rightarrow (\pi, 0)$] is accompanied by the growth of two additional incoherent peaks in the spectral function, from which an anisotropic pseudogap originates.” On general theoretical grounds, in [3, 4] we have shown that the *singular* nature of $n(\mathbf{k})$ at *all* $\mathbf{k} \in \mathcal{S}_F^{(0)}$ (see (iii) below) implies that (see Sec. 10 in [3])

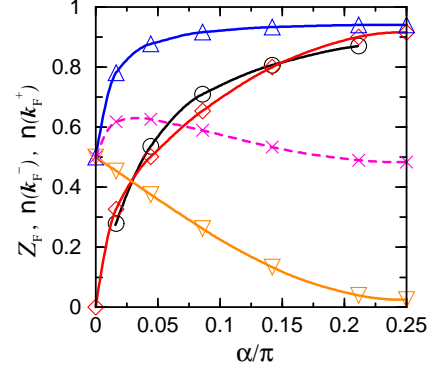


FIG. 1: Z_F as calculated by KK (\circ) and according to the exact expression in the text with $b^- = 0.0912$, $b^+ = -1.4158$ (\diamond), corresponding to $U/t = 2$ and the vH filling associated with $t'/t = 0.1$. $\alpha \equiv \mathbf{k}_F, (\pi, 0)$. The apparent deviation between the two results reflects the isotropy of b^\mp assumed here. Using the above b^\mp , we also present $n(\mathbf{k}_F^-)$ (\triangle) and $n(\mathbf{k}_F^+)$ (∇) obtained from the exact expressions in the text, as well as $[n(\mathbf{k}_F^-) + n(\mathbf{k}_F^+)]/2$ (\times) whose deviation from $1/2$ in the present case is indicative of the underlying metallic state *not* being a Fermi liquid [3] (leaving aside $Z_F = 0$ at $\alpha = 0$).

$Z_F \rightarrow 0$, for $\mathbf{k} \rightarrow \text{PG}$, *must* necessarily be accompanied by at least two *resonant* peaks (to be distinguished from qp peaks) in $A(\mathbf{k}; \varepsilon)$, one *strictly below* and one *strictly above* the Fermi energy ε_F . Here $\mathcal{S}_F^{(0)}$ is the Fermi surface associated with $\varepsilon_{\mathbf{k}}$ (see (iii) below).

(iii) *Fermi surface non-deformation.* — For models involving solely contact-type interaction, we have shown that [3] $\mathcal{S}_F \subseteq \mathcal{S}_F^{(0)}$; PG consists of those points of $\mathcal{S}_F^{(0)}$, if any, which do not belong to \mathcal{S}_F [3, 4]. Interestingly, $\mathcal{S}_F \subseteq \mathcal{S}_F^{(0)}$ turns out to be the working hypothesis for many calculations, amongst which those by KK [1].

We should like to emphasize that the above-mentioned results, cited from [3, 4, 5], are *not* restricted to the weak-coupling limit; the only constraint for the validity of these results is the uniformity of the underlying GSs. We further point out that, *in principle*, depending on the values of d , U/t , t'/t , etc., some other ‘universality classes’ for the uniform metallic GSs of the Hubbard Hamiltonian (explicitly considered in [3, 4]) may/do become viable than that dealt with in this Comment.

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- [5] For cases where $Z_F \neq 0$, some of the pertinent results are more conveniently deduced by the approach in B. Farid, Phil. Mag. **84**, 909 (2004). [cond-mat/0308090]